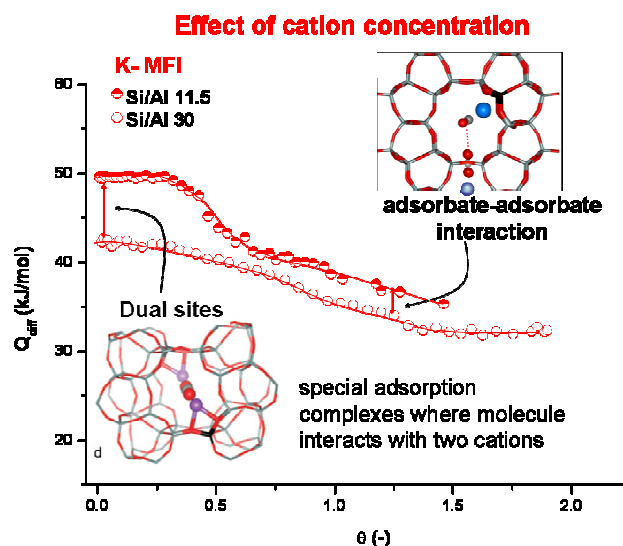
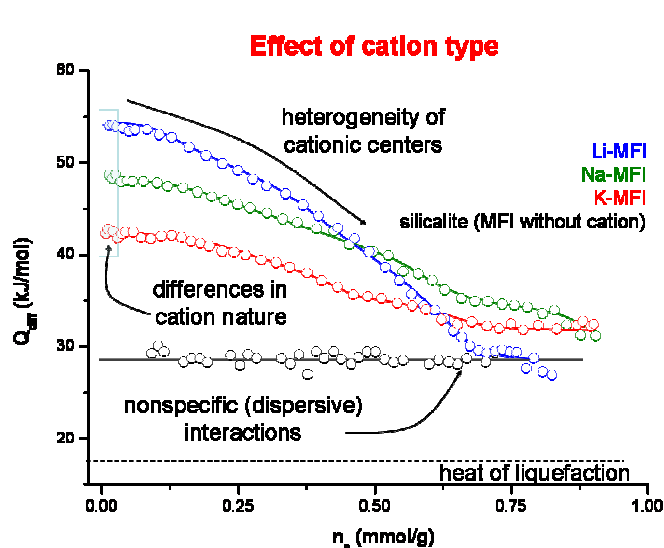


## Study of Adsorption Sites Heterogeneity in Zeolites by Means of Coupled Microcalorimetry with Volumetry

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**Abstract:** Adsorption of  $\text{CO}_2$  as probe molecule on alkali-metal zeolites of MFI structure was investigated by joint volumetry-calorimetry. Consideration was given to the interpretation of the heat evolved when a probe molecule is adsorbed on the surface. In particular, the number and the strength of adsorption sites are discussed as functions of zeolite structure, concentration and nature of extra-framework cation. The adsorption heats ( $q_{iso}$ ) of  $\text{CO}_2$  interaction with alkali-metal cations decrease for MFI zeolite with high Si/Al in the sequence  $\text{Li}^+ > \text{Na}^+ > \text{K}^+$  from 54 kJ/mol to 49 and 43 kJ/mol, respectively. In addition, the adsorption heats are influenced by concentration of Al in the framework. This phenomenon is attributed to formation of bridged  $\text{CO}_2$  adsorption complexes formed between two cations. On the base of quantitative analysis of adsorption processes, presence of geminal adsorption complexes was suggested for adsorption at higher equilibrium pressures.